



Full wwPDB X-ray Structure Validation Report ⓘ

May 19, 2020 – 01:28 pm BST

PDB ID : 6JLG
Title : Crystal Structure of HasAp with Co-9,10,19,20-Tetraphenylporphycene
Authors : Sakakibara, E.; Shisaka, Y.; Onoda, H.; Sugimoto, H.; Shiro, Y.; Watanabe, Y.; Shoji, O.
Deposited on : 2019-03-05
Resolution : 2.50 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.11
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.11

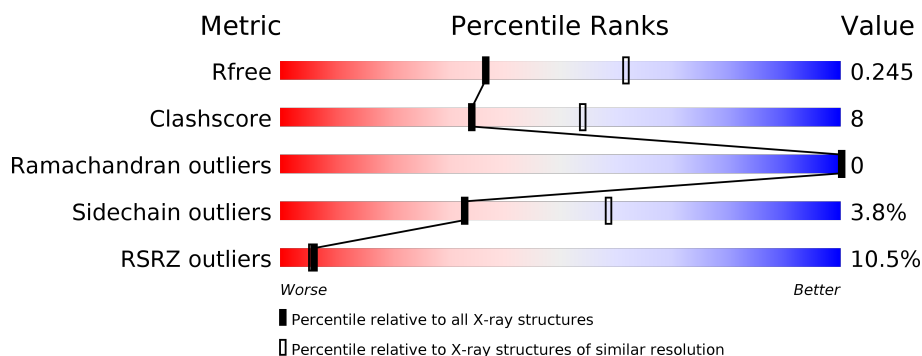
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	4661 (2.50-2.50)
Clashscore	141614	5346 (2.50-2.50)
Ramachandran outliers	138981	5231 (2.50-2.50)
Sidechain outliers	138945	5233 (2.50-2.50)
RSRZ outliers	127900	4559 (2.50-2.50)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	184	<div> <div>11%</div> <div>79%</div> <div>20%</div> <div>.</div> </div>
1	B	184	<div> <div>9%</div> <div>82%</div> <div>16%</div> <div>..</div> </div>
1	C	184	<div> <div>7%</div> <div>82%</div> <div>14%</div> <div>..</div> </div>
1	D	184	<div> <div>13%</div> <div>78%</div> <div>20%</div> <div>.</div> </div>
1	E	184	<div> <div>5%</div> <div>81%</div> <div>17%</div> <div>.</div> </div>
1	F	184	<div> <div>17%</div> <div>72%</div> <div>24%</div> <div>..</div> </div>

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Mol	Chain	Length	Quality of chain
1	G	184	
1	H	184	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	BWU	A	201	X	-	-	-
2	BWU	B	201	X	-	-	-
2	BWU	C	201	X	-	-	-
2	BWU	D	201	X	-	-	-
2	BWU	E	201	X	-	-	-
2	BWU	F	201	X	-	-	-
2	BWU	G	201	X	-	-	-
2	BWU	H	201	X	-	-	-

2 Entry composition

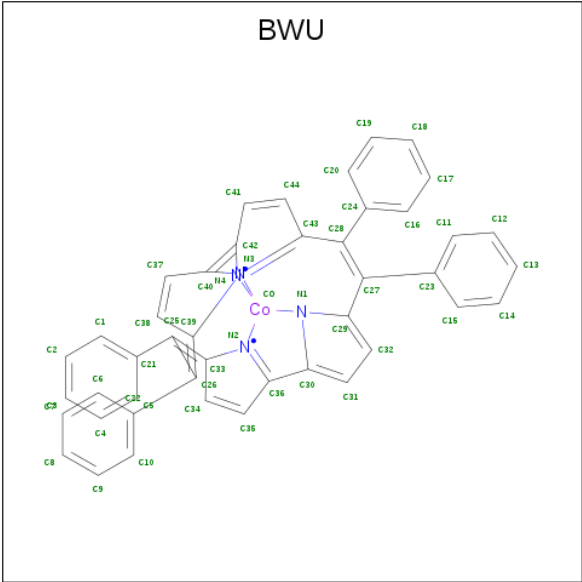
There are 5 unique types of molecules in this entry. The entry contains 11178 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called Heme acquisition protein HasA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	181	Total	C	N	O	S	0	0	0
			1316	825	217	273	1			
1	B	181	Total	C	N	O	S	0	0	0
			1316	825	217	273	1			
1	C	177	Total	C	N	O	S	0	0	0
			1290	812	212	265	1			
1	D	181	Total	C	N	O	S	0	0	0
			1316	825	217	273	1			
1	E	181	Total	C	N	O	S	0	0	0
			1316	825	217	273	1			
1	F	181	Total	C	N	O	S	0	0	0
			1316	825	217	273	1			
1	G	181	Total	C	N	O	S	0	0	0
			1316	825	217	273	1			
1	H	181	Total	C	N	O	S	0	0	0
			1316	825	217	273	1			

- Molecule 2 is Tetraphenylporphycene containing cobalt (three-letter code: BWU) (formula: $C_{44}H_{28}CoN_4$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	Co	N	0	0
			49	44	1	4		
2	B	1	Total	C	Co	N	0	0
			49	44	1	4		
2	C	1	Total	C	Co	N	0	0
			49	44	1	4		
2	D	1	Total	C	Co	N	0	0
			49	44	1	4		
2	E	1	Total	C	Co	N	0	0
			49	44	1	4		
2	F	1	Total	C	Co	N	0	0
			49	44	1	4		
2	G	1	Total	C	Co	N	0	0
			49	44	1	4		
2	H	1	Total	C	Co	N	0	0
			49	44	1	4		

- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			6	3	3		
3	A	1	Total	C	O	0	0
			6	3	3		
3	B	1	Total	C	O	0	0
			6	3	3		
3	B	1	Total	C	O	0	0
			6	3	3		
3	D	1	Total	C	O	0	0
			6	3	3		
3	E	1	Total	C	O	0	0
			6	3	3		
3	G	1	Total	C	O	0	0
			6	3	3		
3	G	1	Total	C	O	0	0
			6	3	3		
3	H	1	Total	C	O	0	0
			6	3	3		

- Molecule 4 is PHOSPHATE ION (three-letter code: PO4) (formula: O₄P).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	E	1	Total	O	P	0	0
			5	4	1		
4	G	1	Total	O	P	0	0
			5	4	1		

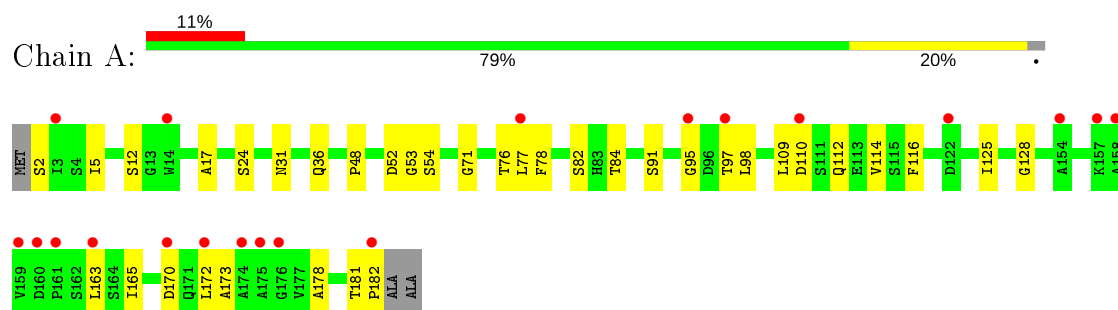
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	32	Total	O	0	0
			32	32		
5	B	36	Total	O	0	0
			36	36		
5	C	18	Total	O	0	0
			18	18		
5	D	17	Total	O	0	0
			17	17		
5	E	47	Total	O	0	0
			47	47		
5	F	11	Total	O	0	0
			11	11		
5	G	41	Total	O	0	0
			41	41		
5	H	18	Total	O	0	0
			18	18		

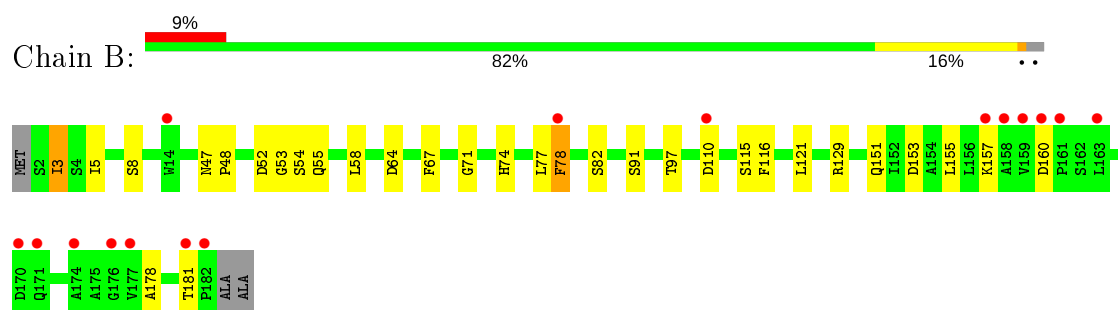
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

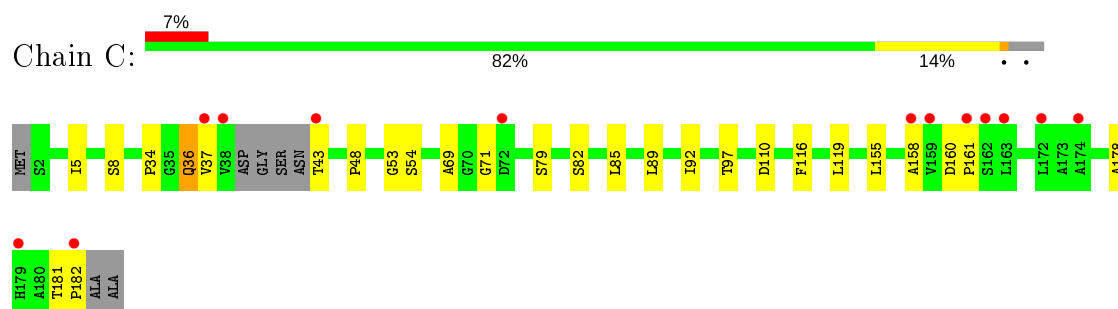
- Molecule 1: Heme acquisition protein HasA



- Molecule 1: Heme acquisition protein HasA

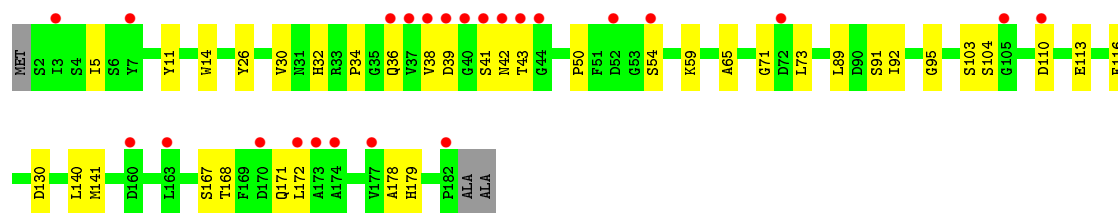


- Molecule 1: Heme acquisition protein HasA

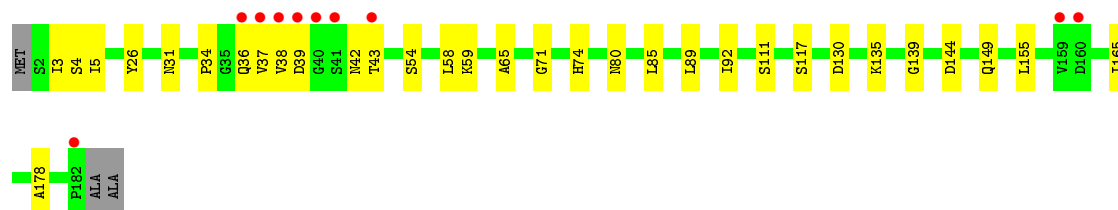
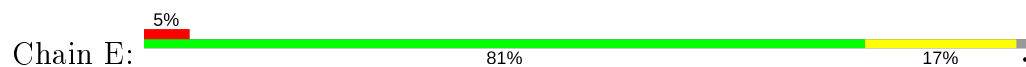


- Molecule 1: Heme acquisition protein HasA

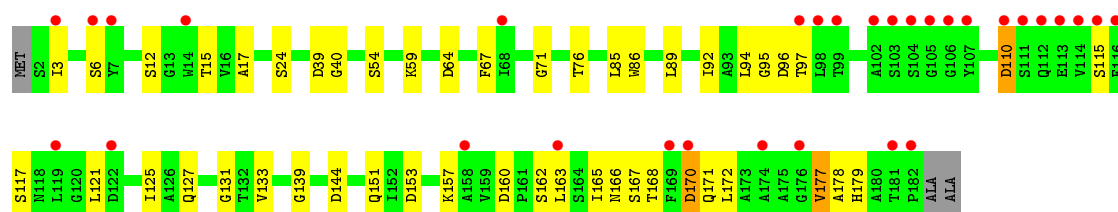




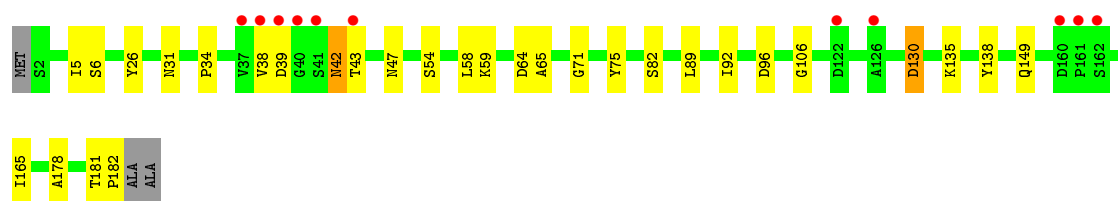
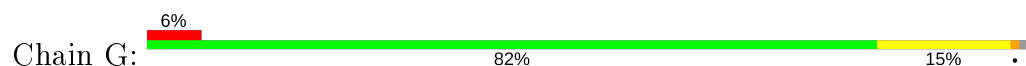
• Molecule 1: Heme acquisition protein HasA



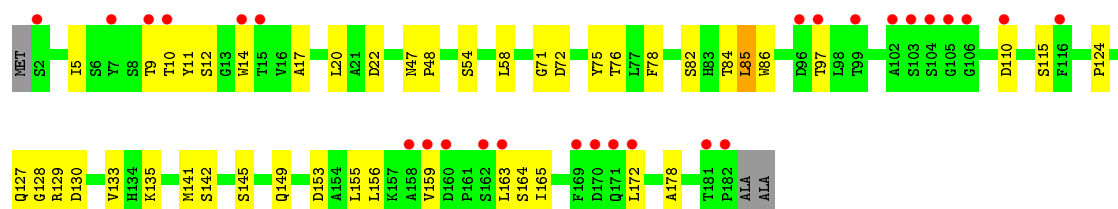
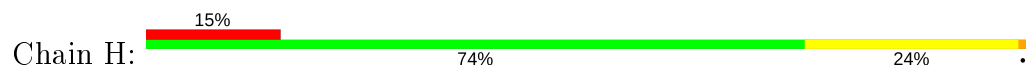
• Molecule 1: Heme acquisition protein HasA



• Molecule 1: Heme acquisition protein HasA



• Molecule 1: Heme acquisition protein HasA



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	237.58Å 92.22Å 93.19Å 90.00° 100.96° 90.00°	Depositor
Resolution (Å)	45.79 – 2.50 45.75 – 2.50	Depositor EDS
% Data completeness (in resolution range)	99.7 (45.79-2.50) 99.7 (45.75-2.50)	Depositor EDS
R_{merge}	0.25	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.78 (at 2.51Å)	Xtriage
Refinement program	REFMAC 5.8.0238	Depositor
R, R_{free}	0.213 , 0.241 0.217 , 0.245	Depositor DCC
R_{free} test set	3400 reflections (4.97%)	wwPDB-VP
Wilson B-factor (Å ²)	42.5	Xtriage
Anisotropy	0.436	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 48.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	11178	wwPDB-VP
Average B, all atoms (Å ²)	54.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 19.00% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: GOL, PO4, BWU

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.64	0/1348	0.67	0/1840
1	B	0.64	0/1348	0.67	0/1840
1	C	0.64	0/1321	0.66	0/1802
1	D	0.65	0/1348	0.66	0/1840
1	E	0.65	0/1348	0.67	0/1840
1	F	0.64	0/1348	0.65	0/1840
1	G	0.65	0/1348	0.67	0/1840
1	H	0.64	0/1348	0.66	0/1840
All	All	0.65	0/10757	0.67	0/14682

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1316	0	1229	23	0
1	B	1316	0	1229	24	0
1	C	1290	0	1209	14	0
1	D	1316	0	1228	26	0
1	E	1316	0	1228	20	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	F	1316	0	1229	32	0
1	G	1316	0	1228	17	0
1	H	1316	0	1229	25	0
2	A	49	0	0	0	0
2	B	49	0	0	0	0
2	C	49	0	0	0	0
2	D	49	0	0	1	0
2	E	49	0	0	0	0
2	F	49	0	0	0	0
2	G	49	0	0	1	0
2	H	49	0	0	2	0
3	A	12	0	16	1	0
3	B	12	0	16	0	0
3	D	6	0	8	2	0
3	E	6	0	8	0	0
3	G	12	0	16	1	0
3	H	6	0	8	0	0
4	E	5	0	0	0	0
4	G	5	0	0	1	0
5	A	32	0	0	0	0
5	B	36	0	0	0	0
5	C	18	0	0	0	0
5	D	17	0	0	0	0
5	E	47	0	0	0	0
5	F	11	0	0	0	0
5	G	41	0	0	0	0
5	H	18	0	0	0	0
All	All	11178	0	9881	174	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 8.

All (174) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:3:ILE:HG12	1:B:155:LEU:HB3	1.36	1.04
1:B:3:ILE:H	1:B:3:ILE:HD12	1.38	0.86
1:G:42:ASN:O	1:G:42:ASN:ND2	2.09	0.85
1:A:2:SER:OG	3:A:202:GOL:H2	1.77	0.84
1:B:3:ILE:CG1	1:B:155:LEU:HB3	2.07	0.83
1:A:112:GLN:NE2	1:A:114:VAL:C	2.33	0.82

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:160:ASP:OD1	1:C:161:PRO:HD2	1.82	0.80
1:H:156:LEU:HB3	1:H:163:LEU:HD12	1.65	0.78
1:F:168:THR:HB	1:F:171:GLN:HG3	1.66	0.77
1:G:43:THR:O	1:G:58:LEU:HD12	1.85	0.77
1:A:112:GLN:HE22	1:A:114:VAL:C	1.87	0.76
1:H:129:ARG:O	1:H:135:LYS:HE3	1.90	0.72
1:A:95:GLY:HA3	1:A:109:LEU:HD13	1.73	0.71
1:H:156:LEU:O	1:H:159:VAL:HG12	1.91	0.70
1:D:32:HIS:CE1	1:D:39:ASP:HB3	2.27	0.69
1:F:97:THR:HB	1:F:110:ASP:HB3	1.75	0.68
1:A:97:THR:HB	1:A:110:ASP:HB3	1.77	0.65
1:A:112:GLN:NE2	1:A:114:VAL:O	2.30	0.65
1:F:6:SER:HB2	1:F:115:SER:OG	1.97	0.64
1:D:34:PRO:HG2	1:E:34:PRO:HD2	1.79	0.64
1:A:181:THR:HB	1:A:182:PRO:HD2	1.81	0.63
1:F:3:ILE:HD12	1:F:117:SER:O	2.00	0.62
1:D:38:VAL:HG21	1:D:42:ASN:ND2	2.15	0.61
1:E:38:VAL:HG22	1:E:39:ASP:H	1.65	0.61
1:D:36:GLN:HG2	1:D:38:VAL:O	2.00	0.61
1:A:54:SER:HA	1:A:71:GLY:O	2.00	0.60
1:F:139:GLY:C	1:F:144:ASP:O	2.40	0.60
1:F:153:ASP:OD1	1:F:165:ILE:N	2.35	0.59
1:B:78:PHE:CE2	3:D:202:GOL:H12	2.37	0.59
1:H:153:ASP:HB2	1:H:165:ILE:HD11	1.84	0.59
1:H:20:LEU:HD12	1:H:149:GLN:HG2	1.85	0.58
1:H:163:LEU:HD11	1:H:172:LEU:HD21	1.86	0.58
1:C:181:THR:HG23	1:C:182:PRO:HD2	1.84	0.57
1:C:48:PRO:HD2	1:C:53:GLY:HA2	1.87	0.56
1:B:78:PHE:HZ	1:D:50:PRO:HB3	1.69	0.56
1:F:15:THR:HG23	1:F:166:ASN:HA	1.87	0.56
1:A:31:ASN:HB3	1:A:36:GLN:OE1	2.05	0.56
1:G:149:GLN:HG3	1:G:165:ILE:HG12	1.87	0.56
1:E:89:LEU:HD21	1:E:92:ILE:HD11	1.88	0.55
1:B:3:ILE:CD1	1:B:3:ILE:H	2.08	0.55
1:E:43:THR:O	1:E:59:LYS:N	2.36	0.55
1:F:172:LEU:O	1:F:177:VAL:HG12	2.06	0.55
1:E:38:VAL:HG22	1:E:39:ASP:N	2.22	0.54
1:B:77:LEU:O	1:B:82:SER:HA	2.06	0.54
1:D:168:THR:O	1:D:172:LEU:HD22	2.08	0.54
1:E:26:TYR:CD2	1:E:65:ALA:HB2	2.43	0.54
1:D:89:LEU:HD21	1:D:92:ILE:HD11	1.90	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:97:THR:HB	1:H:110:ASP:HB3	1.90	0.54
1:A:76:THR:HG21	1:A:125:ILE:CD1	2.37	0.53
1:A:181:THR:O	1:A:182:PRO:C	2.45	0.53
1:B:78:PHE:CZ	3:D:202:GOL:H12	2.44	0.53
1:C:116:PHE:HB3	1:C:119:LEU:HD11	1.89	0.53
1:D:95:GLY:HA2	1:D:113:GLU:HG3	1.91	0.53
1:C:36:GLN:HG2	1:C:36:GLN:O	2.08	0.53
1:F:139:GLY:CA	1:F:144:ASP:O	2.57	0.52
1:H:11:TYR:HA	1:H:14:TRP:CD1	2.44	0.51
1:C:160:ASP:OD1	1:C:161:PRO:CD	2.54	0.51
1:E:3:ILE:CG2	1:E:155:LEU:HD13	2.41	0.51
1:E:43:THR:HG23	1:E:58:LEU:HD11	1.92	0.51
1:B:52:ASP:OD2	1:B:74:HIS:ND1	2.35	0.50
1:C:54:SER:HA	1:C:71:GLY:O	2.11	0.50
1:F:89:LEU:HD21	1:F:92:ILE:HD11	1.92	0.50
1:D:32:HIS:HA	1:D:36:GLN:HE22	1.76	0.50
1:B:153:ASP:OD2	1:B:157:LYS:HE2	2.11	0.50
1:H:124:PRO:HG2	1:H:127:GLN:HG3	1.92	0.50
1:H:54:SER:HA	1:H:71:GLY:O	2.10	0.50
1:D:26:TYR:CD2	1:D:65:ALA:HB2	2.47	0.50
1:G:59:LYS:HE2	1:G:64:ASP:HA	1.93	0.49
1:B:97:THR:HB	1:B:110:ASP:HB3	1.94	0.49
1:B:129:ARG:HD3	1:G:26:TYR:CG	2.46	0.49
1:H:145:SER:O	1:H:149:GLN:HG3	2.12	0.49
1:B:5:ILE:O	1:B:178:ALA:HA	2.13	0.49
1:E:36:GLN:CG	1:E:36:GLN:O	2.61	0.49
1:H:76:THR:HG23	1:H:86:TRP:HE1	1.78	0.49
1:B:54:SER:HA	1:B:71:GLY:O	2.13	0.49
1:F:76:THR:HG21	1:F:125:ILE:CD1	2.43	0.49
1:A:48:PRO:HD2	1:A:53:GLY:HA2	1.94	0.48
1:F:157:LYS:HA	1:F:160:ASP:O	2.13	0.48
1:C:69:ALA:HB1	1:C:89:LEU:HD11	1.95	0.48
1:E:5:ILE:O	1:E:178:ALA:HA	2.14	0.48
1:F:54:SER:HA	1:F:71:GLY:O	2.14	0.48
1:B:91:SER:HA	1:B:116:PHE:O	2.12	0.48
1:G:130:ASP:O	1:G:135:LYS:NZ	2.46	0.48
1:G:42:ASN:OD1	1:H:78:PHE:CE1	2.66	0.48
1:F:127:GLN:HE21	1:F:131:GLY:HA2	1.77	0.48
1:D:92:ILE:HD13	1:D:140:LEU:HD11	1.95	0.47
1:G:82:SER:OG	4:G:202:PO4:O1	2.28	0.47
1:G:26:TYR:CD2	1:G:65:ALA:HB2	2.49	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:163:LEU:HD13	1:H:164:SER:N	2.30	0.47
1:D:54:SER:HA	1:D:71:GLY:O	2.15	0.47
1:H:17:ALA:N	1:H:165:ILE:HG22	2.30	0.47
1:C:155:LEU:O	1:C:158:ALA:HB3	2.15	0.47
1:A:95:GLY:CA	1:A:109:LEU:HD13	2.43	0.46
1:D:41:SER:HB3	1:D:59:LYS:HB3	1.97	0.46
1:G:5:ILE:O	1:G:178:ALA:HA	2.15	0.46
1:G:89:LEU:HD21	1:G:92:ILE:HD11	1.96	0.46
1:D:168:THR:O	1:D:172:LEU:CD2	2.64	0.46
1:F:76:THR:HG21	1:F:125:ILE:HD12	1.97	0.46
1:D:30:VAL:HG22	1:D:141:MET:O	2.15	0.46
1:E:149:GLN:CG	1:E:165:ILE:HG12	2.46	0.46
1:F:67:PHE:HB2	1:F:94:LEU:HD23	1.98	0.46
1:A:97:THR:O	1:A:98:LEU:C	2.54	0.46
1:E:54:SER:HA	1:E:71:GLY:O	2.16	0.46
1:D:167:SER:HB2	1:D:172:LEU:HD21	1.98	0.46
1:B:3:ILE:HG12	1:B:155:LEU:CB	2.25	0.45
1:F:160:ASP:OD1	1:F:162:SER:N	2.43	0.45
1:A:76:THR:HG21	1:A:125:ILE:HD13	1.98	0.45
1:F:95:GLY:C	1:F:96:ASP:OD1	2.55	0.45
1:D:34:PRO:CG	1:E:34:PRO:HD2	2.46	0.45
1:H:9:THR:O	1:H:12:SER:OG	2.28	0.45
1:E:4:SER:OG	1:E:117:SER:HB3	2.17	0.45
1:B:121:LEU:HD11	1:B:151:GLN:HG2	1.98	0.44
1:F:167:SER:HB2	1:F:172:LEU:HD21	1.99	0.44
1:F:15:THR:HG22	1:F:17:ALA:H	1.82	0.44
1:F:76:THR:HG23	1:F:86:TRP:HE1	1.82	0.44
1:B:3:ILE:N	1:B:3:ILE:HD12	2.19	0.44
1:C:5:ILE:O	1:C:178:ALA:HA	2.18	0.44
1:H:72:ASP:N	1:H:72:ASP:OD1	2.51	0.44
1:B:160:ASP:OD1	1:B:160:ASP:N	2.48	0.44
1:F:15:THR:HG22	1:F:17:ALA:N	2.32	0.44
1:A:163:LEU:HD22	1:A:172:LEU:HD23	1.99	0.44
1:C:34:PRO:HG2	1:G:34:PRO:HD2	2.00	0.44
1:D:11:TYR:CE2	1:D:113:GLU:HB3	2.53	0.44
1:E:36:GLN:O	1:E:36:GLN:HG3	2.18	0.44
1:G:138:TYR:HE2	3:G:204:GOL:H31	1.83	0.43
1:B:48:PRO:HD2	1:B:53:GLY:HA2	1.99	0.43
1:C:97:THR:HB	1:C:110:ASP:HB3	1.99	0.43
1:G:75:TYR:CZ	2:G:201:BWU:N3	2.86	0.43
1:B:3:ILE:CD1	1:B:155:LEU:HB3	2.47	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:47:ASN:ND2	1:G:106:GLY:HA2	2.34	0.43
1:F:160:ASP:HB3	1:F:163:LEU:HD12	2.01	0.43
1:H:85:LEU:HD23	1:H:133:VAL:HG12	2.00	0.43
1:F:85:LEU:HD23	1:F:133:VAL:HG12	2.00	0.43
1:A:5:ILE:O	1:A:178:ALA:HA	2.18	0.43
1:E:130:ASP:O	1:E:135:LYS:HE3	2.17	0.43
1:D:11:TYR:OH	1:H:130:ASP:OD2	2.30	0.43
1:E:74:HIS:O	1:E:85:LEU:HA	2.19	0.42
1:H:5:ILE:O	1:H:178:ALA:HA	2.20	0.42
1:B:8:SER:HA	1:B:181:THR:O	2.19	0.42
1:E:139:GLY:CA	1:E:144:ASP:O	2.67	0.42
1:F:121:LEU:HD11	1:F:151:GLN:HG2	2.01	0.42
1:D:168:THR:HG23	1:D:171:GLN:H	1.83	0.42
1:A:84:THR:HG21	1:A:128:GLY:HA2	2.01	0.42
1:D:32:HIS:CD2	2:D:201:BWU:N4	2.87	0.42
1:D:38:VAL:HG21	1:D:42:ASN:HD21	1.84	0.42
1:D:130:ASP:OD2	1:F:59:LYS:NZ	2.52	0.42
1:H:47:ASN:HA	1:H:48:PRO:HA	1.90	0.42
1:C:79:SER:O	1:C:82:SER:HA	2.20	0.41
1:F:39:ASP:OD1	1:F:40:GLY:N	2.50	0.41
1:H:75:TYR:CE1	2:H:201:BWU:C40	3.03	0.41
1:F:178:ALA:C	1:F:179:HIS:HD2	2.23	0.41
1:F:95:GLY:O	1:F:96:ASP:OD1	2.38	0.41
1:G:181:THR:HG23	1:G:182:PRO:HD2	2.02	0.41
1:D:11:TYR:HA	1:D:14:TRP:CD1	2.56	0.41
1:A:12:SER:HB3	1:A:170:ASP:OD2	2.21	0.41
1:E:43:THR:HG23	1:E:58:LEU:CD1	2.51	0.41
1:A:77:LEU:O	1:A:82:SER:HA	2.21	0.41
1:B:47:ASN:HB3	1:B:55:GLN:HB2	2.03	0.41
1:B:58:LEU:HB3	1:B:67:PHE:CE2	2.55	0.41
1:E:165:ILE:HD12	1:E:165:ILE:HA	1.88	0.41
1:D:91:SER:HA	1:D:116:PHE:O	2.21	0.40
1:G:54:SER:HA	1:G:71:GLY:O	2.21	0.40
1:F:12:SER:HB2	1:F:170:ASP:OD1	2.21	0.40
1:H:84:THR:HG21	1:H:128:GLY:HA2	2.03	0.40
1:A:173:ALA:HA	1:A:178:ALA:O	2.20	0.40
1:D:5:ILE:O	1:D:178:ALA:HA	2.21	0.40
1:F:17:ALA:HB2	1:F:165:ILE:HG23	2.03	0.40
2:H:201:BWU:C10	2:H:201:BWU:C21	2.97	0.40
1:H:20:LEU:HA	1:H:20:LEU:HD23	1.88	0.40
1:A:17:ALA:HB2	1:A:165:ILE:HG23	2.03	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:91:SER:HA	1:A:116:PHE:O	2.21	0.40
1:C:89:LEU:HD21	1:C:92:ILE:HD11	2.03	0.40
1:F:172:LEU:HA	1:F:172:LEU:HD13	1.89	0.40
1:H:58:LEU:HD22	1:H:141:MET:HG2	2.04	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	179/184 (97%)	172 (96%)	7 (4%)	0	100	100
1	B	179/184 (97%)	171 (96%)	8 (4%)	0	100	100
1	C	173/184 (94%)	168 (97%)	5 (3%)	0	100	100
1	D	179/184 (97%)	173 (97%)	6 (3%)	0	100	100
1	E	179/184 (97%)	170 (95%)	9 (5%)	0	100	100
1	F	179/184 (97%)	177 (99%)	2 (1%)	0	100	100
1	G	179/184 (97%)	172 (96%)	7 (4%)	0	100	100
1	H	179/184 (97%)	176 (98%)	3 (2%)	0	100	100
All	All	1426/1472 (97%)	1379 (97%)	47 (3%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was

analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	138/139 (99%)	135 (98%)	3 (2%)	52	77
1	B	138/139 (99%)	134 (97%)	4 (3%)	42	69
1	C	135/139 (97%)	130 (96%)	5 (4%)	34	60
1	D	138/139 (99%)	132 (96%)	6 (4%)	29	53
1	E	138/139 (99%)	133 (96%)	5 (4%)	35	61
1	F	138/139 (99%)	133 (96%)	5 (4%)	35	61
1	G	138/139 (99%)	131 (95%)	7 (5%)	24	45
1	H	138/139 (99%)	131 (95%)	7 (5%)	24	45
All	All	1101/1112 (99%)	1059 (96%)	42 (4%)	33	58

All (42) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	24	SER
1	A	52	ASP
1	A	78	PHE
1	B	3	ILE
1	B	64	ASP
1	B	78	PHE
1	B	115	SER
1	C	8	SER
1	C	36	GLN
1	C	37	VAL
1	C	43	THR
1	C	85	LEU
1	D	43	THR
1	D	73	LEU
1	D	103	SER
1	D	104	SER
1	D	110	ASP
1	D	179	HIS
1	E	31	ASN
1	E	37	VAL
1	E	42	ASN
1	E	80	ASN
1	E	111	SER
1	F	24	SER
1	F	64	ASP

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Mol	Chain	Res	Type
1	F	110	ASP
1	F	170	ASP
1	F	177	VAL
1	G	6	SER
1	G	31	ASN
1	G	38	VAL
1	G	39	ASP
1	G	42	ASN
1	G	96	ASP
1	G	130	ASP
1	H	10	THR
1	H	22	ASP
1	H	82	SER
1	H	85	LEU
1	H	115	SER
1	H	142	SER
1	H	155	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	112	GLN
1	D	36	GLN
1	D	42	ASN
1	E	42	ASN
1	F	112	GLN
1	F	127	GLN
1	F	179	HIS
1	G	74	HIS
1	G	166	ASN
1	H	127	GLN
1	H	171	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

19 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z > 2$	Counts	RMSZ	# $ Z > 2$
3	GOL	B	202	-	5,5,5	0.12	0	5,5,5	0.28	0
3	GOL	G	204	-	5,5,5	0.15	0	5,5,5	0.44	0
3	GOL	D	202	-	5,5,5	0.11	0	5,5,5	0.29	0
2	BWU	D	201	1	49,60,60	2.86	4 (8%)	51,94,94	1.95	9 (17%)
3	GOL	H	202	-	5,5,5	0.10	0	5,5,5	0.27	0
3	GOL	A	203	-	5,5,5	0.10	0	5,5,5	0.25	0
4	PO4	G	202	-	4,4,4	0.67	0	6,6,6	0.43	0
3	GOL	E	203	-	5,5,5	0.12	0	5,5,5	0.31	0
3	GOL	B	203	-	5,5,5	0.14	0	5,5,5	0.32	0
3	GOL	G	203	-	5,5,5	0.14	0	5,5,5	0.35	0
3	GOL	A	202	-	5,5,5	0.11	0	5,5,5	0.29	0
4	PO4	E	202	-	4,4,4	0.71	0	6,6,6	0.46	0
2	BWU	G	201	1	49,60,60	2.88	4 (8%)	51,94,94	1.96	8 (15%)
2	BWU	E	201	1	49,60,60	2.79	4 (8%)	51,94,94	1.95	9 (17%)
2	BWU	F	201	1	49,60,60	3.09	4 (8%)	51,94,94	1.95	9 (17%)
2	BWU	C	201	1	49,60,60	2.96	4 (8%)	51,94,94	1.93	8 (15%)
2	BWU	A	201	1	49,60,60	2.97	3 (6%)	51,94,94	1.95	9 (17%)
2	BWU	B	201	1	49,60,60	3.03	4 (8%)	51,94,94	2.03	9 (17%)
2	BWU	H	201	1	49,60,60	3.04	4 (8%)	51,94,94	1.95	9 (17%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	GOL	B	202	-	-	1/4/4/4	-
3	GOL	G	204	-	-	0/4/4/4	-
3	GOL	D	202	-	-	0/4/4/4	-
2	BWU	D	201	1	1/1/5/5	0/16/72/72	0/4/12/12
3	GOL	A	203	-	-	2/4/4/4	-
3	GOL	H	202	-	-	0/4/4/4	-
3	GOL	E	203	-	-	0/4/4/4	-
3	GOL	B	203	-	-	1/4/4/4	-
3	GOL	G	203	-	-	0/4/4/4	-
3	GOL	A	202	-	-	1/4/4/4	-
2	BWU	G	201	1	1/1/5/5	0/16/72/72	0/4/12/12
2	BWU	E	201	1	1/1/5/5	0/16/72/72	0/4/12/12
2	BWU	F	201	1	1/1/5/5	0/16/72/72	0/4/12/12
2	BWU	C	201	1	1/1/5/5	0/16/72/72	0/4/12/12
2	BWU	A	201	1	1/1/5/5	0/16/72/72	0/4/12/12
2	BWU	B	201	1	1/1/5/5	0/16/72/72	0/4/12/12
2	BWU	H	201	1	1/1/5/5	0/16/72/72	0/4/12/12

All (31) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	F	201	BWU	C25-C26	-20.61	1.33	1.49
2	H	201	BWU	C25-C26	-20.33	1.33	1.49
2	B	201	BWU	C25-C26	-20.21	1.34	1.49
2	A	201	BWU	C25-C26	-19.84	1.34	1.49
2	C	201	BWU	C25-C26	-19.81	1.34	1.49
2	D	201	BWU	C25-C26	-19.05	1.34	1.49
2	G	201	BWU	C25-C26	-18.92	1.34	1.49
2	E	201	BWU	C25-C26	-18.49	1.35	1.49
2	G	201	BWU	C36-C30	-4.13	1.37	1.47
2	E	201	BWU	C36-C30	-3.80	1.38	1.47
2	F	201	BWU	C36-C30	-3.58	1.39	1.47
2	H	201	BWU	C36-C30	-3.54	1.39	1.47
2	A	201	BWU	C36-C30	-3.51	1.39	1.47
2	B	201	BWU	C36-C30	-3.50	1.39	1.47
2	D	201	BWU	C36-C30	-3.37	1.39	1.47
2	C	201	BWU	C36-C30	-3.34	1.39	1.47
2	G	201	BWU	C36-N2	-2.40	1.33	1.36
2	G	201	BWU	C42-N4	-2.40	1.33	1.38
2	F	201	BWU	C42-N4	-2.31	1.34	1.38
2	H	201	BWU	C42-N4	-2.31	1.34	1.38
2	E	201	BWU	C42-N4	-2.28	1.34	1.38
2	A	201	BWU	C42-N4	-2.27	1.34	1.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	201	BWU	C36-N2	-2.26	1.33	1.36
2	D	201	BWU	C42-N4	-2.22	1.34	1.38
2	B	201	BWU	C42-N4	-2.19	1.34	1.38
2	E	201	BWU	C36-N2	-2.18	1.33	1.36
2	D	201	BWU	C36-N2	-2.18	1.33	1.36
2	C	201	BWU	C42-N4	-2.16	1.34	1.38
2	C	201	BWU	C36-N2	-2.14	1.33	1.36
2	H	201	BWU	C36-N2	-2.13	1.33	1.36
2	F	201	BWU	C36-N2	-2.08	1.34	1.36

All (70) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	201	BWU	C36-C30-N1	6.74	124.18	115.08
2	H	201	BWU	C36-C30-N1	6.70	124.11	115.08
2	D	201	BWU	C36-C30-N1	6.64	124.03	115.08
2	E	201	BWU	C36-C30-N1	6.42	123.74	115.08
2	A	201	BWU	C36-C30-N1	6.41	123.72	115.08
2	G	201	BWU	C36-C30-N1	6.29	123.57	115.08
2	B	201	BWU	C36-C30-N1	6.25	123.50	115.08
2	C	201	BWU	C30-C36-N2	6.23	123.48	115.08
2	C	201	BWU	C36-C30-N1	6.18	123.41	115.08
2	F	201	BWU	C30-C36-N2	6.02	123.21	115.08
2	B	201	BWU	C30-C36-N2	6.02	123.19	115.08
2	H	201	BWU	C30-C36-N2	5.95	123.10	115.08
2	A	201	BWU	C30-C36-N2	5.89	123.03	115.08
2	G	201	BWU	C30-C36-N2	5.78	122.88	115.08
2	B	201	BWU	C32-C31-C30	-5.77	103.83	107.61
2	D	201	BWU	C30-C36-N2	5.73	122.81	115.08
2	E	201	BWU	C34-C35-C36	-5.42	104.06	107.61
2	E	201	BWU	C30-C36-N2	5.40	122.37	115.08
2	B	201	BWU	C34-C35-C36	-5.26	104.16	107.61
2	G	201	BWU	C34-C35-C36	-5.24	104.17	107.61
2	F	201	BWU	C34-C35-C36	-5.03	104.31	107.61
2	A	201	BWU	C32-C31-C30	-5.00	104.33	107.61
2	H	201	BWU	C34-C35-C36	-4.95	104.36	107.61
2	G	201	BWU	C32-C31-C30	-4.84	104.43	107.61
2	C	201	BWU	C34-C35-C36	-4.79	104.47	107.61
2	A	201	BWU	C34-C35-C36	-4.79	104.47	107.61
2	D	201	BWU	C34-C35-C36	-4.70	104.53	107.61
2	H	201	BWU	C32-C31-C30	-4.69	104.54	107.61
2	C	201	BWU	C32-C31-C30	-4.57	104.62	107.61

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	201	BWU	C32-C31-C30	-4.42	104.71	107.61
2	F	201	BWU	C32-C31-C30	-4.32	104.78	107.61
2	D	201	BWU	C44-C41-C42	-4.26	103.99	107.37
2	D	201	BWU	C32-C31-C30	-3.98	105.00	107.61
2	B	201	BWU	C44-C41-C42	-3.97	104.22	107.37
2	C	201	BWU	C38-C37-C40	-3.86	104.31	107.37
2	C	201	BWU	C44-C41-C42	-3.83	104.33	107.37
2	E	201	BWU	C38-C37-C40	-3.69	104.44	107.37
2	F	201	BWU	C38-C37-C40	-3.65	104.47	107.37
2	D	201	BWU	C38-C37-C40	-3.65	104.47	107.37
2	B	201	BWU	C31-C30-C36	-3.60	123.30	132.03
2	H	201	BWU	C31-C30-C36	-3.52	123.48	132.03
2	G	201	BWU	C38-C37-C40	-3.50	104.59	107.37
2	A	201	BWU	C31-C30-C36	-3.49	123.55	132.03
2	F	201	BWU	C31-C30-C36	-3.47	123.61	132.03
2	G	201	BWU	C31-C30-C36	-3.43	123.69	132.03
2	B	201	BWU	C35-C36-C30	-3.41	123.75	132.03
2	C	201	BWU	C35-C36-C30	-3.37	123.84	132.03
2	E	201	BWU	C31-C30-C36	-3.37	123.84	132.03
2	G	201	BWU	C35-C36-C30	-3.37	123.85	132.03
2	F	201	BWU	C35-C36-C30	-3.35	123.90	132.03
2	D	201	BWU	C31-C30-C36	-3.34	123.92	132.03
2	F	201	BWU	C44-C41-C42	-3.32	104.73	107.37
2	E	201	BWU	C44-C41-C42	-3.28	104.77	107.37
2	A	201	BWU	C38-C37-C40	-3.28	104.77	107.37
2	H	201	BWU	C35-C36-C30	-3.27	124.09	132.03
2	A	201	BWU	C35-C36-C30	-3.26	124.11	132.03
2	C	201	BWU	C31-C30-C36	-3.25	124.13	132.03
2	H	201	BWU	C38-C37-C40	-3.24	104.80	107.37
2	A	201	BWU	C44-C41-C42	-3.24	104.80	107.37
2	E	201	BWU	C35-C36-C30	-3.21	124.24	132.03
2	D	201	BWU	C35-C36-C30	-3.18	124.31	132.03
2	H	201	BWU	C44-C41-C42	-3.15	104.87	107.37
2	B	201	BWU	C38-C37-C40	-2.76	105.18	107.37
2	B	201	BWU	C24-C28-C27	-2.64	116.12	120.27
2	A	201	BWU	C24-C28-C27	-2.60	116.17	120.27
2	G	201	BWU	C44-C41-C42	-2.59	105.31	107.37
2	D	201	BWU	C24-C28-C27	-2.53	116.29	120.27
2	H	201	BWU	C24-C28-C27	-2.42	116.46	120.27
2	F	201	BWU	C24-C28-C27	-2.41	116.48	120.27
2	E	201	BWU	C24-C28-C27	-2.05	117.04	120.27

All (8) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	D	201	BWU	N3
2	G	201	BWU	N3
2	E	201	BWU	N3
2	F	201	BWU	N3
2	C	201	BWU	N3
2	A	201	BWU	N3
2	B	201	BWU	N3
2	H	201	BWU	N3

All (5) torsion outliers are listed below:

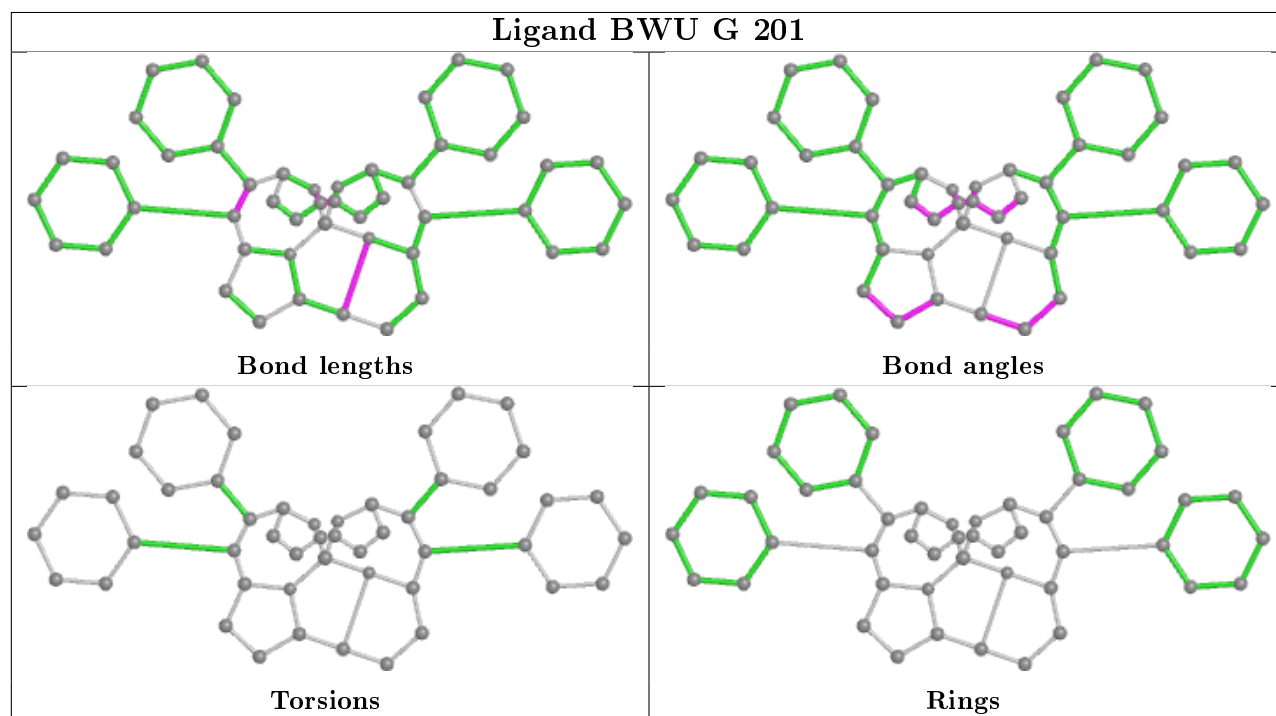
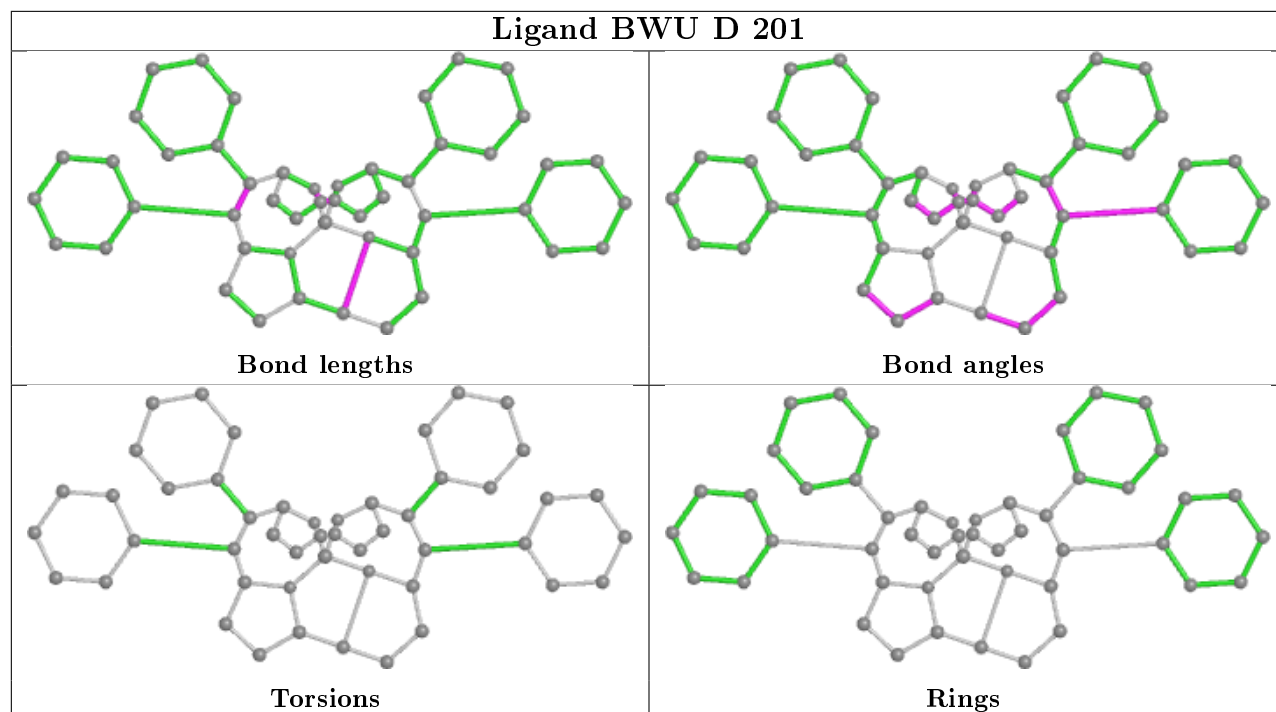
Mol	Chain	Res	Type	Atoms
3	A	203	GOL	O2-C2-C3-O3
3	B	203	GOL	O1-C1-C2-O2
3	A	202	GOL	O2-C2-C3-O3
3	B	202	GOL	O1-C1-C2-C3
3	A	203	GOL	C1-C2-C3-O3

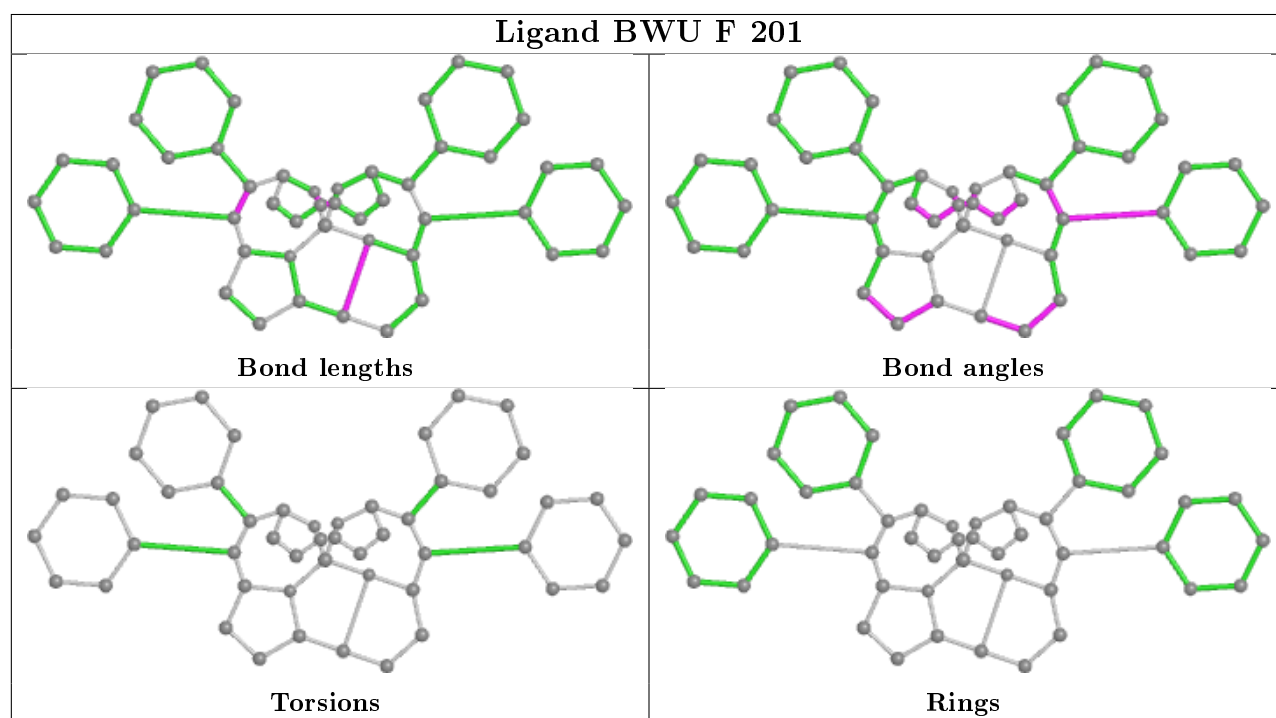
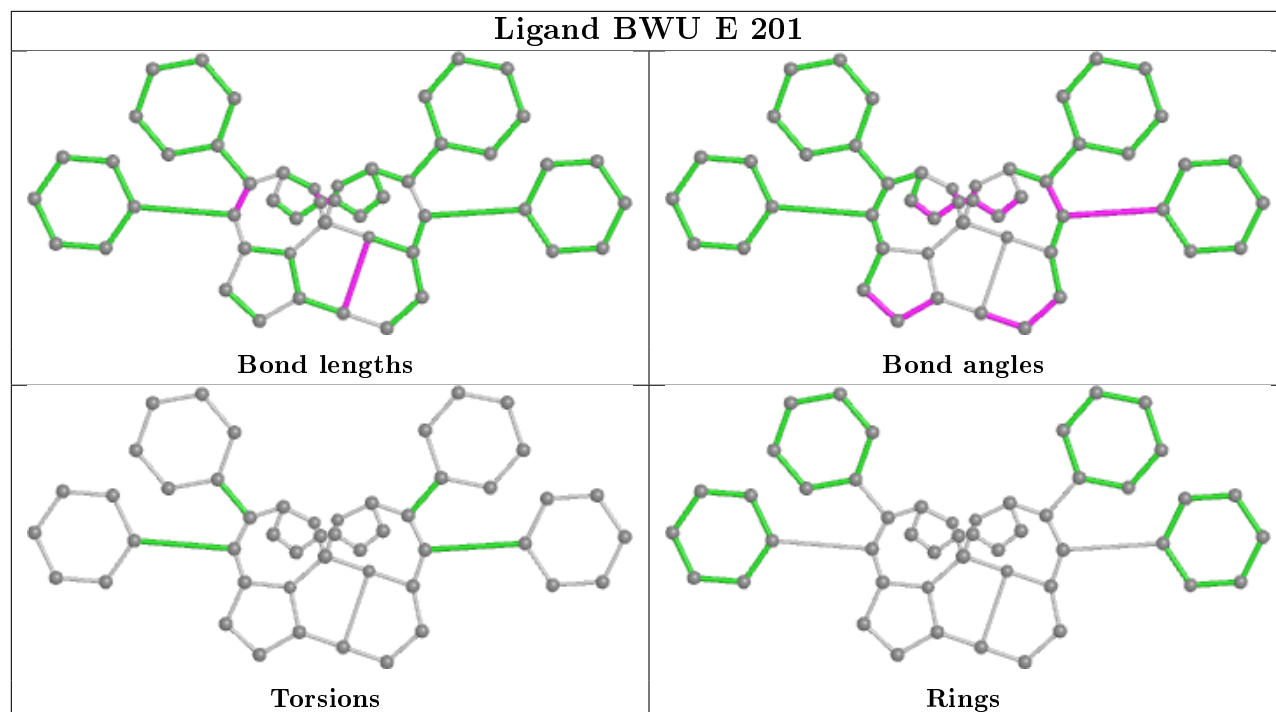
There are no ring outliers.

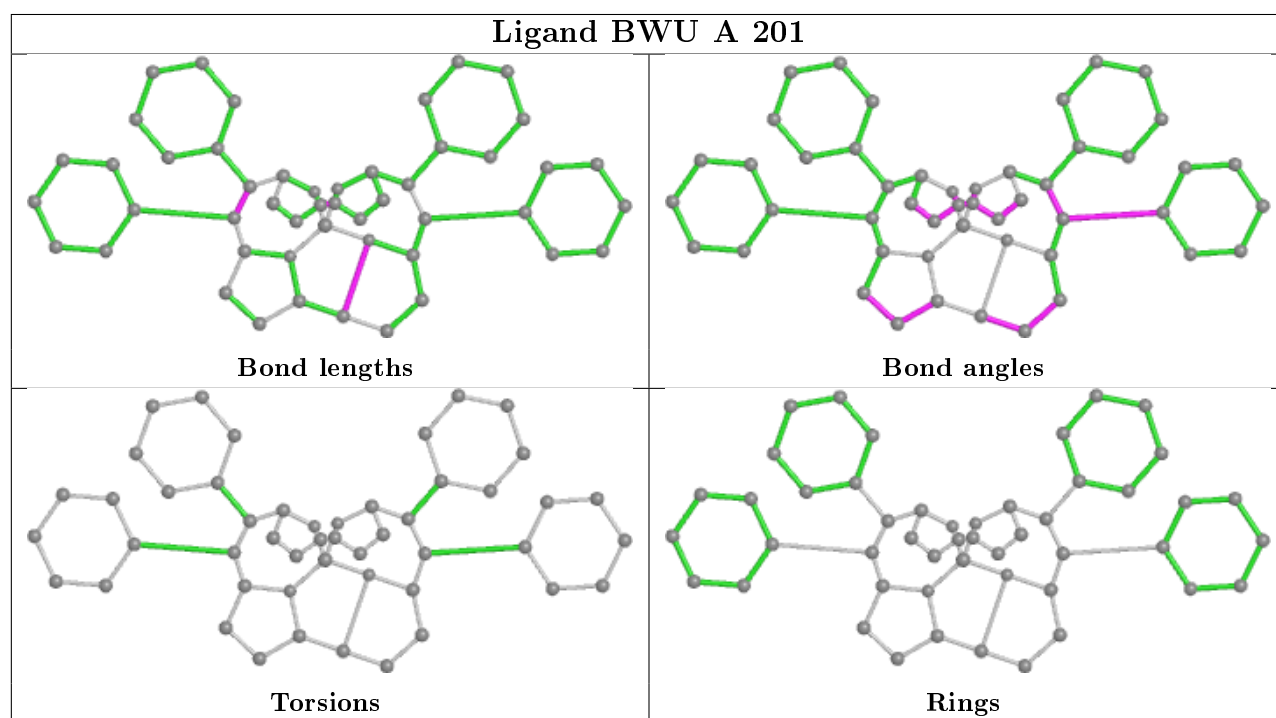
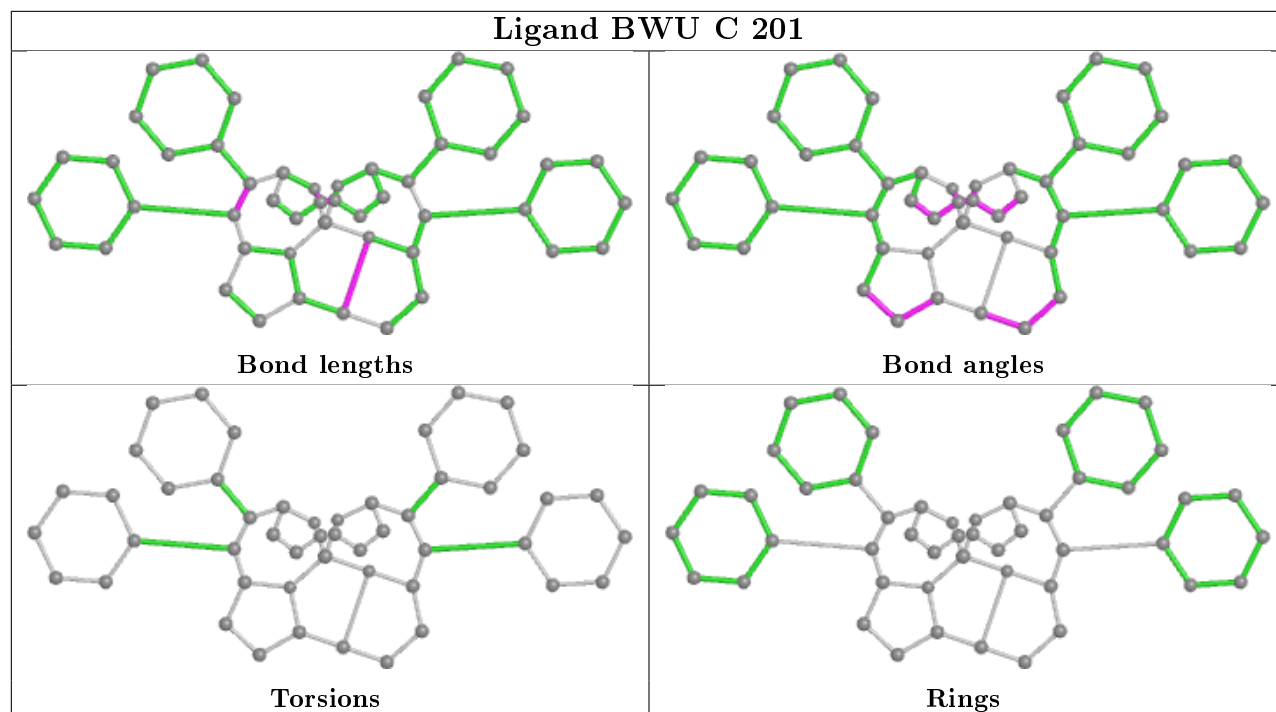
7 monomers are involved in 9 short contacts:

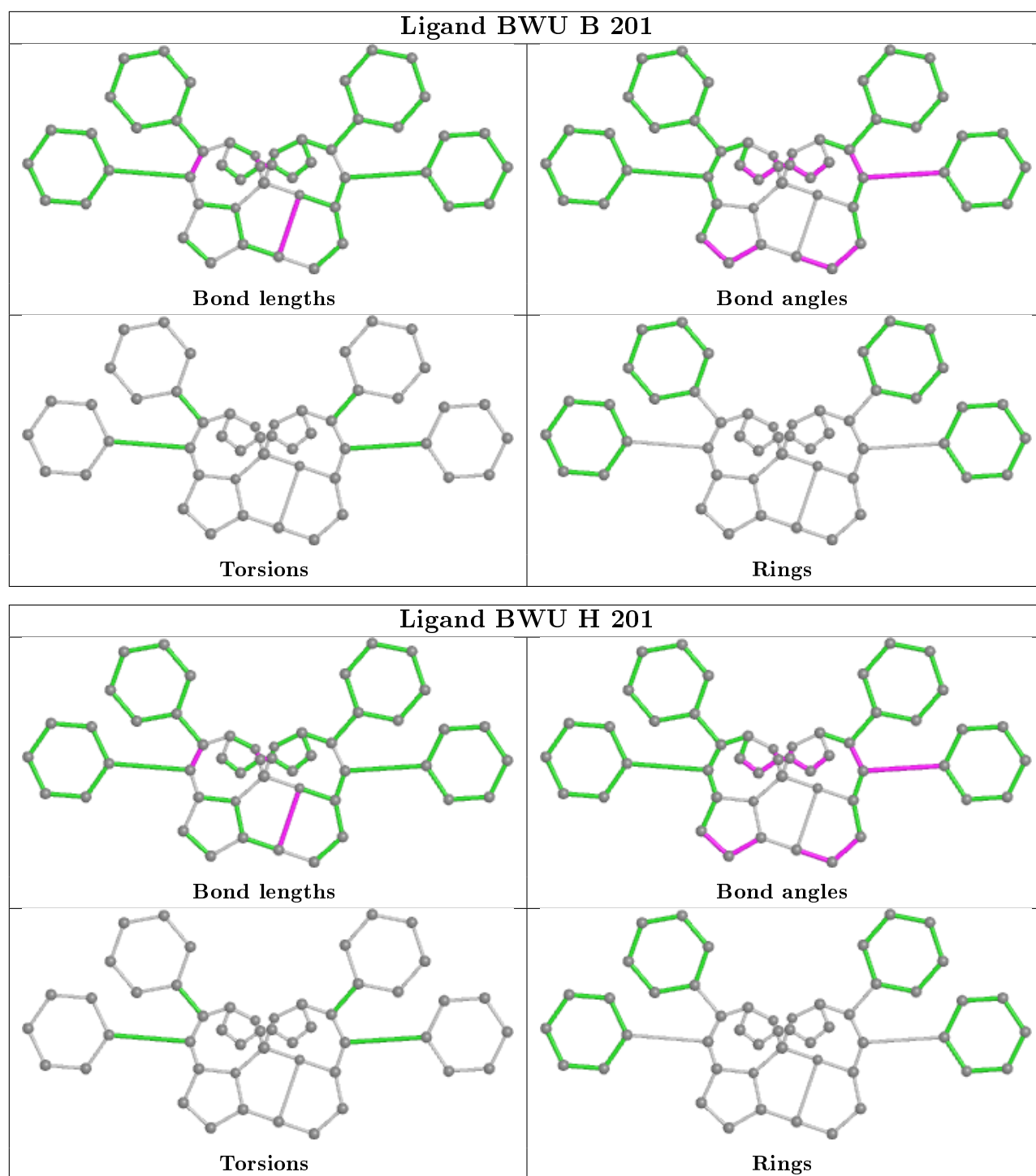
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	G	204	GOL	1	0
3	D	202	GOL	2	0
2	D	201	BWU	1	0
4	G	202	PO4	1	0
3	A	202	GOL	1	0
2	G	201	BWU	1	0
2	H	201	BWU	2	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.









5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	181/184 (98%)	0.61	20 (11%) 5 5	28, 45, 81, 117	0
1	B	181/184 (98%)	0.69	16 (8%) 10 10	28, 44, 84, 112	0
1	C	177/184 (96%)	0.53	13 (7%) 15 15	34, 55, 86, 95	0
1	D	181/184 (98%)	0.77	24 (13%) 3 2	31, 55, 92, 123	0
1	E	181/184 (98%)	0.40	10 (5%) 25 26	27, 40, 63, 132	0
1	F	181/184 (98%)	0.89	31 (17%) 1 1	39, 66, 110, 126	0
1	G	181/184 (98%)	0.45	11 (6%) 21 22	28, 41, 65, 136	0
1	H	181/184 (98%)	0.83	27 (14%) 2 2	37, 60, 93, 112	0
All	All	1444/1472 (98%)	0.65	152 (10%) 6 6	27, 50, 94, 136	0

All (152) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	H	163	LEU	7.0
1	C	163	LEU	7.0
1	D	38	VAL	5.8
1	G	40	GLY	5.7
1	D	40	GLY	5.7
1	D	39	ASP	5.6
1	B	170	ASP	5.5
1	B	160	ASP	5.4
1	E	37	VAL	5.4
1	H	102	ALA	5.3
1	F	102	ALA	5.1
1	H	14	TRP	5.1
1	B	182	PRO	5.0
1	B	159	VAL	5.0
1	G	39	ASP	4.8
1	G	37	VAL	4.7

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Mol	Chain	Res	Type	RSRZ
1	B	158	ALA	4.6
1	G	38	VAL	4.6
1	D	41	SER	4.6
1	A	159	VAL	4.4
1	F	106	GLY	4.3
1	C	37	VAL	4.3
1	B	176	GLY	4.2
1	F	110	ASP	4.2
1	D	172	LEU	4.2
1	A	176	GLY	4.2
1	H	181	THR	4.2
1	F	97	THR	4.1
1	G	41	SER	4.1
1	B	177	VAL	4.1
1	D	37	VAL	4.0
1	H	104	SER	3.8
1	B	181	THR	3.8
1	H	162	SER	3.8
1	A	3	ILE	3.8
1	H	103	SER	3.7
1	A	160	ASP	3.7
1	H	159	VAL	3.7
1	D	44	GLY	3.7
1	F	6	SER	3.7
1	F	182	PRO	3.6
1	F	3	ILE	3.6
1	H	170	ASP	3.6
1	A	95	GLY	3.6
1	A	161	PRO	3.6
1	D	42	ASN	3.5
1	A	170	ASP	3.5
1	E	41	SER	3.5
1	A	122	ASP	3.4
1	D	72	ASP	3.4
1	F	99	THR	3.4
1	G	160	ASP	3.4
1	C	38	VAL	3.4
1	A	157	LYS	3.4
1	C	182	PRO	3.3
1	D	163	LEU	3.3
1	D	43	THR	3.3
1	B	163	LEU	3.3

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Mol	Chain	Res	Type	RSRZ
1	E	38	VAL	3.3
1	B	174	ALA	3.2
1	F	170	ASP	3.1
1	F	98	LEU	3.1
1	H	105	GLY	3.1
1	E	160	ASP	3.1
1	A	110	ASP	3.0
1	D	170	ASP	3.0
1	F	116	PHE	3.0
1	A	158	ALA	3.0
1	H	169	PHE	3.0
1	B	161	PRO	3.0
1	H	106	GLY	2.9
1	E	39	ASP	2.9
1	D	174	ALA	2.9
1	D	36	GLN	2.9
1	D	182	PRO	2.9
1	A	175	ALA	2.8
1	D	160	ASP	2.8
1	C	43	THR	2.8
1	H	116	PHE	2.8
1	F	176	GLY	2.8
1	F	14	TRP	2.8
1	F	105	GLY	2.8
1	F	163	LEU	2.8
1	C	158	ALA	2.8
1	D	52	ASP	2.7
1	C	179	HIS	2.7
1	E	40	GLY	2.7
1	E	36	GLN	2.6
1	C	172	LEU	2.6
1	F	169	PHE	2.6
1	H	10	THR	2.6
1	F	107	TYR	2.6
1	H	96	ASP	2.6
1	D	54	SER	2.6
1	F	111	SER	2.6
1	F	68	ILE	2.5
1	E	43	THR	2.5
1	H	160	ASP	2.5
1	G	43	THR	2.5
1	F	104	SER	2.5

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Mol	Chain	Res	Type	RSRZ
1	H	9	THR	2.5
1	B	110	ASP	2.5
1	B	14	TRP	2.4
1	D	110	ASP	2.4
1	H	172	LEU	2.4
1	G	126	ALA	2.4
1	C	72	ASP	2.4
1	F	103	SER	2.4
1	B	78	PHE	2.4
1	E	182	PRO	2.4
1	H	110	ASP	2.4
1	H	15	THR	2.4
1	A	163	LEU	2.4
1	A	14	TRP	2.3
1	F	7	TYR	2.3
1	D	177	VAL	2.3
1	H	158	ALA	2.3
1	G	161	PRO	2.3
1	G	162	SER	2.3
1	F	181	THR	2.3
1	C	161	PRO	2.3
1	H	97	THR	2.3
1	F	113	GLU	2.3
1	H	2	SER	2.2
1	C	174	ALA	2.2
1	A	77	LEU	2.2
1	B	157	LYS	2.2
1	H	99	THR	2.2
1	H	182	PRO	2.2
1	F	114	VAL	2.2
1	H	7	TYR	2.2
1	E	159	VAL	2.2
1	F	158	ALA	2.2
1	C	159	VAL	2.1
1	A	174	ALA	2.1
1	F	174	ALA	2.1
1	C	162	SER	2.1
1	D	3	ILE	2.1
1	G	122	ASP	2.1
1	A	182	PRO	2.1
1	A	154	ALA	2.1
1	A	172	LEU	2.1

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Mol	Chain	Res	Type	RSRZ
1	F	119	LEU	2.1
1	D	173	ALA	2.0
1	A	97	THR	2.0
1	D	7	TYR	2.0
1	B	171	GLN	2.0
1	F	112	GLN	2.0
1	F	115	SER	2.0
1	D	105	GLY	2.0
1	F	122	ASP	2.0
1	H	171	GLN	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

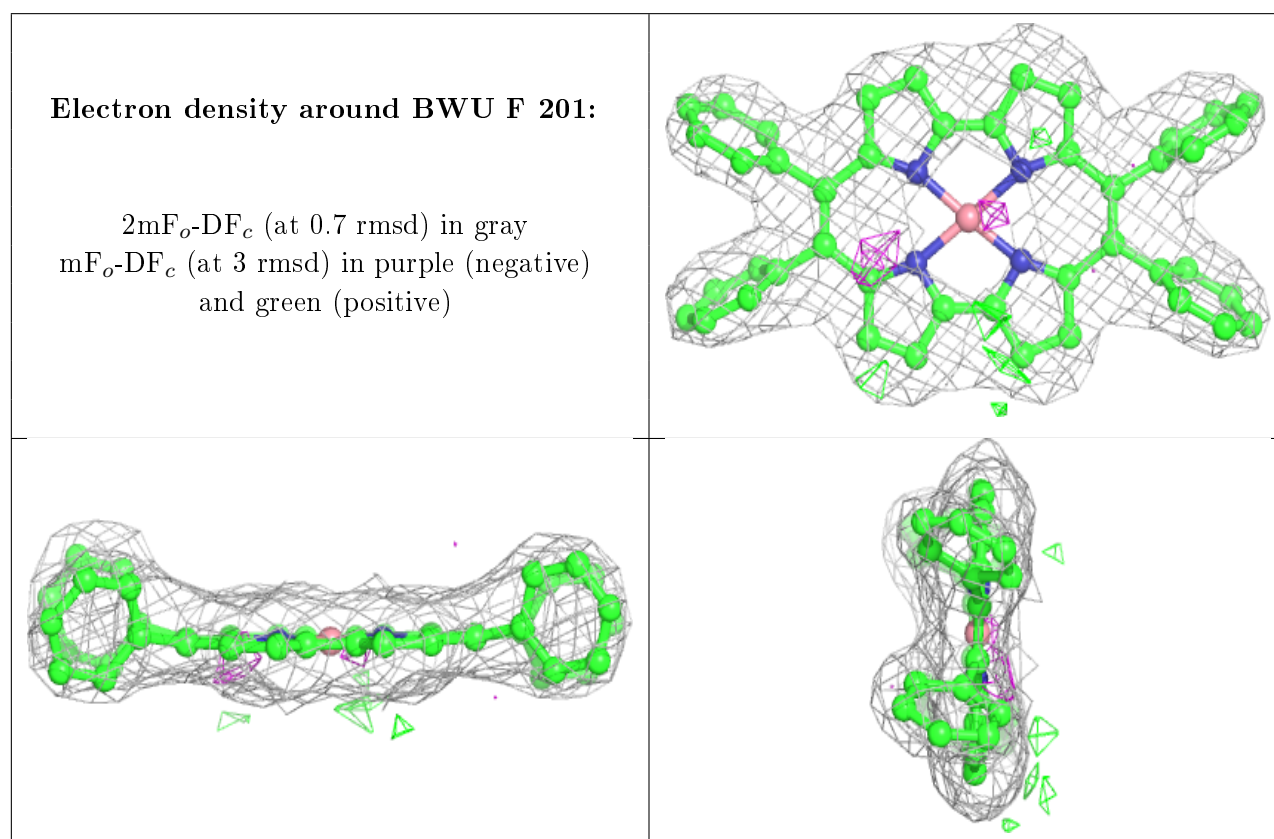
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	GOL	B	202	6/6	0.74	0.18	45,57,58,58	0
3	GOL	E	203	6/6	0.74	0.22	50,55,56,57	0
3	GOL	G	203	6/6	0.74	0.28	42,52,56,56	0
3	GOL	G	204	6/6	0.75	0.27	39,47,48,50	0
3	GOL	A	202	6/6	0.76	0.22	51,54,55,56	0
3	GOL	B	203	6/6	0.83	0.23	40,48,50,52	0
3	GOL	D	202	6/6	0.89	0.19	41,47,50,50	0
3	GOL	A	203	6/6	0.89	0.15	48,49,52,53	0
3	GOL	H	202	6/6	0.90	0.23	51,55,55,58	0
2	BWU	F	201	49/49	0.95	0.17	37,40,49,51	0
4	PO4	G	202	5/5	0.96	0.21	35,41,46,53	0
2	BWU	A	201	49/49	0.96	0.17	27,31,34,35	0
2	BWU	B	201	49/49	0.96	0.17	26,30,34,36	0

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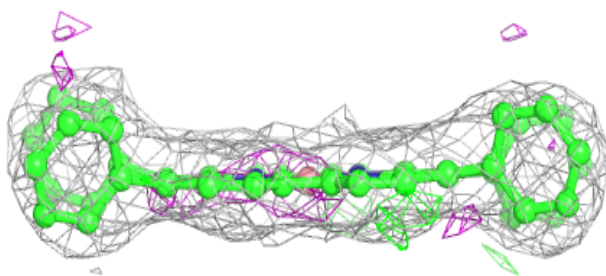
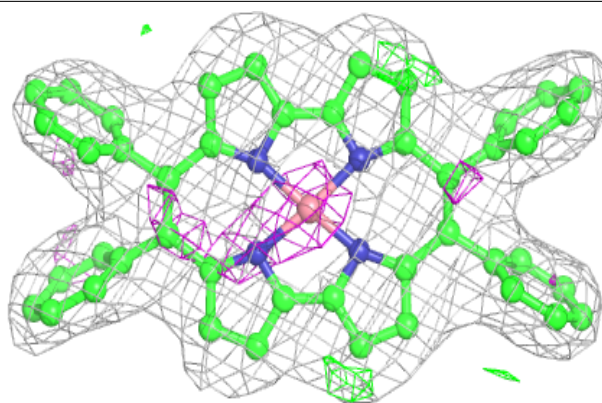
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	BWU	H	201	49/49	0.96	0.16	36,40,45,46	0
2	BWU	C	201	49/49	0.97	0.16	30,35,50,56	0
2	BWU	G	201	49/49	0.97	0.16	26,31,39,41	0
2	BWU	E	201	49/49	0.97	0.17	24,28,38,39	0
2	BWU	D	201	49/49	0.97	0.16	28,31,51,57	0
4	PO4	E	202	5/5	0.98	0.21	34,41,45,55	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

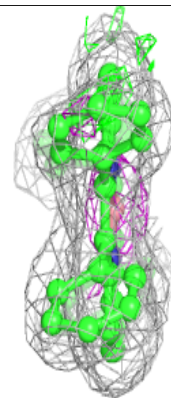
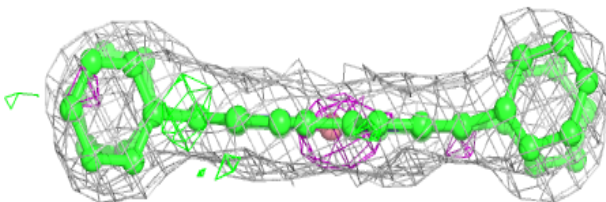
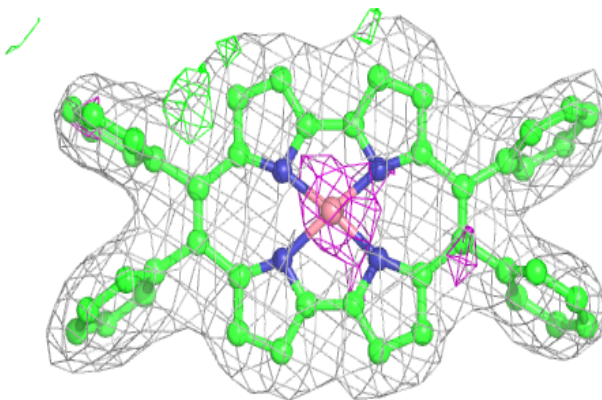


Electron density around BWU A 201:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

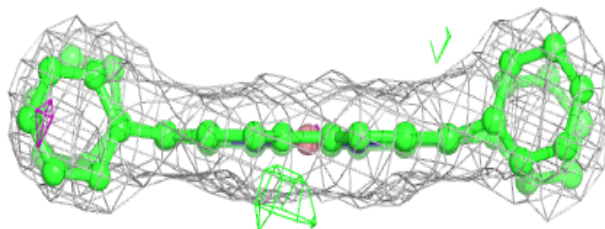
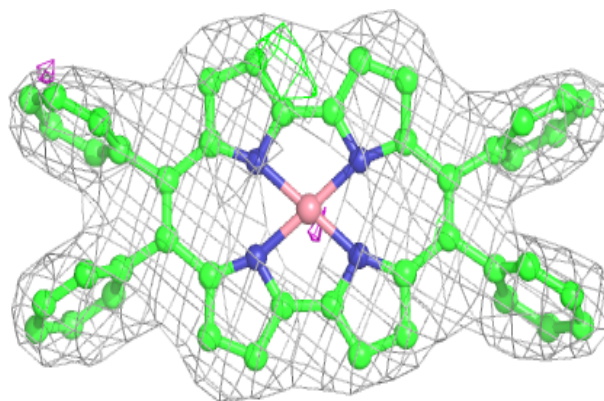
**Electron density around BWU B 201:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

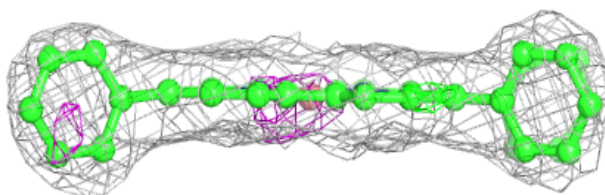
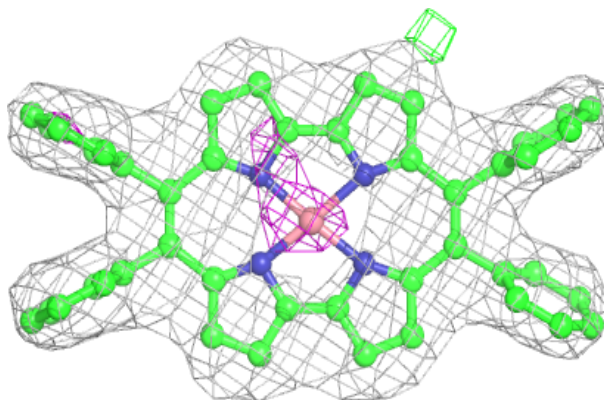


Electron density around BWU H 201:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

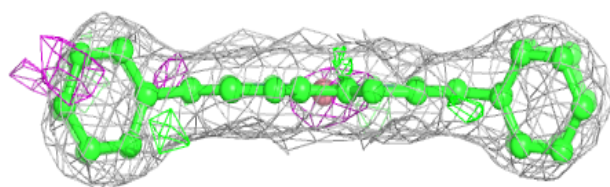
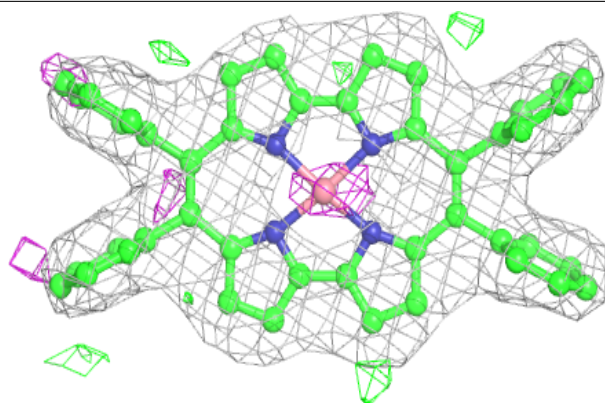
**Electron density around BWU C 201:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

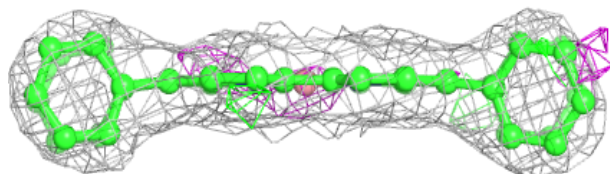
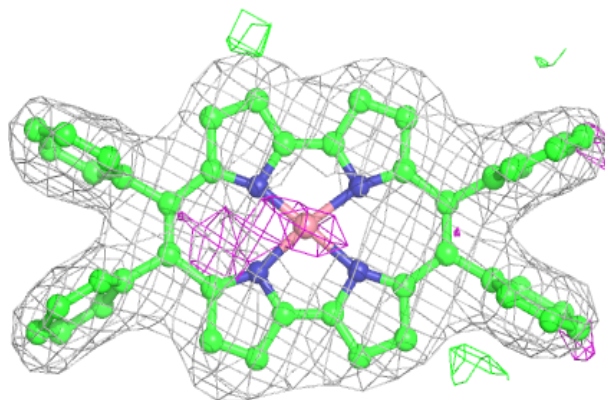


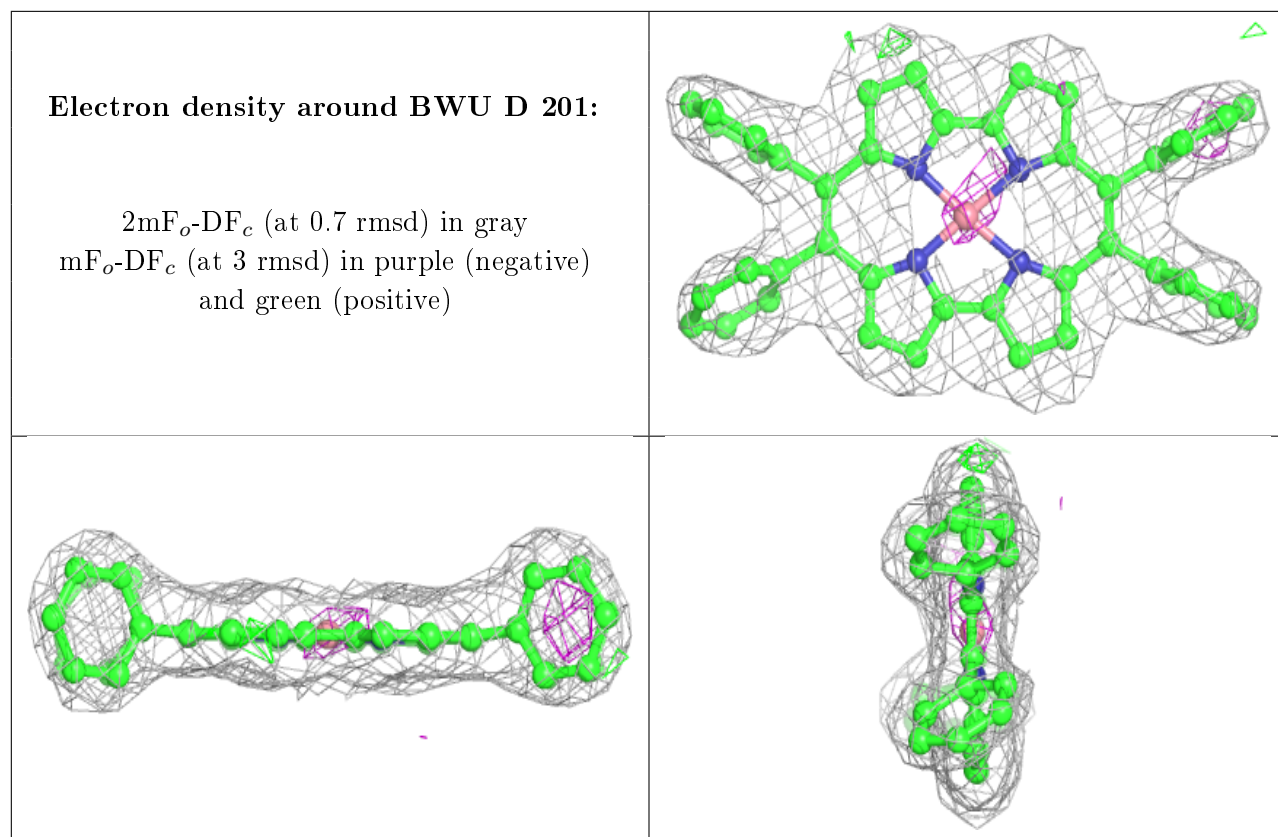
Electron density around BWU G 201:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

**Electron density around BWU E 201:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)





6.5 Other polymers [i](#)

There are no such residues in this entry.